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Minimizing Statistical Bias with Queries

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Abstract

I describe an exploration criterion that attempts to minimize the error of a learner by minimizing its estimated squared bias. I describe experiments with locally-weighted regression on two simple kinematics problems, and observe that this “bias-only” approach outperforms the more common “variance-only” exploration approach, even in the presence of noise.

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1 Introduction

In recent years, there has been an explosion of interest “active” machine learning systems. These are learning systems that make queries, or perform experiments to gather data that are expected to maximize performance. When compared with “passive” learning systems, which accept given, or randomly drawn data, active learners have demonstrated significant decreases in the amount of data required to achieve equivalent performance. In industrial applications, where each experiment may take days to perform and cost thousands of dollars, a method for optimally selecting these points would offer enormous savings in time and money.

An active learning system will typically attempt to select data that will minimize its predictive error. The error of a learner can be decomposed into bias and variance terms. Most research in selecting optimal actions or queries has assumed that the learner is approximately unbiased, and that to minimize learner error, variance is the only thing to minimize (a few examples include Fedorov [1972], MacKay [1992], Cohn [1994; 1995], Paass [1995]). In practice, however, there are very few problems for which we have unbiased learners. Frequently, bias constitutes a large portion of a learner’s error; if the learner is deterministic and the data are noise-free, then bias is the *only* source of error.¹

In this paper I describe an algorithm which selects actions/queries designed to minimize the bias of a locally weighted regression-based learner. Empirically, “variance-minimizing” strategies which ignore bias seem to perform well, even in cases where, strictly speaking, there is no variance to minimize. In the tasks considered in this paper, the bias-minimizing strategy consistently outperforms variance minimization, even in the presence of noise.

1.1 Bias and variance

Let us begin by defining $P(x, y)$ to be the unknown joint distribution over x and y , and $P(x)$ to be the known marginal distribution of x (commonly called the *input distribution*). We denote the learner’s output on input x , given training set \mathcal{D} as $\hat{y}(x; \mathcal{D})$. We can then write the expected error of the learner as

$$\int_x E \left[(\hat{y}(x; \mathcal{D}) - y(x))^2 | x \right] P(x) dx, \quad (1)$$

where $E[\cdot]$ denotes the expectation over P and over training sets \mathcal{D} . The expectation inside the integral may be decomposed as follows (Geman et al., 1992):

$$E \left[(\hat{y}(x; \mathcal{D}) - y(x))^2 | x \right] = \quad (2)$$

$$\begin{aligned} & E \left[(y(x) - E[y|x])^2 \right] \\ & + (E_{\mathcal{D}}[y(x; \mathcal{D})] - E[y|x])^2 \\ & + E_{\mathcal{D}} \left[(\hat{y}(x; \mathcal{D}) - E_{\mathcal{D}}[\hat{y}(x; \mathcal{D})])^2 \right] \end{aligned} \quad (3)$$

¹The bias term here is a *statistical* bias, which is distinct from the *inductive* bias discussed in some machine learning research. See Dietterich and Kong [1995] for a discussion of the relationship between the two.

where $E_{\mathcal{D}}[\cdot]$ denotes the expectation over training sets. The first term in Equation 2 is the variance of y given x – it is the *noise* in the distribution, and does not depend on our learner or how the training data are chosen. The second term is the learner’s *squared bias*, and the third is its *variance*; these last two terms comprise the expected squared error of the learner with respect to the regression function $E[y|x]$.

Most research in active learning assumes that the second term of Equation 2 is approximately zero, that is, that the learner is *unbiased*. If this is the case, then one may concentrate on selecting data so as to minimize the variance of the learner. Although this “all-variance” approach is optimal when the learner is unbiased, truly unbiased learners are rare. Even when the learner’s representation class is able to match the target function exactly, bias is generally introduced by the learning algorithm and learning parameters. From the Bayesian perspective, a learner is only unbiased if its priors are *exactly* correct.

The optimal choice of query would, of course, minimize *both* bias and variance, but I leave that for future work. For the purposes of this paper, I will only be concerned with selecting queries that are expected to minimize learner bias. This approach is justified in cases where noise is believed to be only a small component of the learner’s error. If the learner is deterministic and there is no noise, then strictly speaking, there *is* no error due to variance — all the error must be due to learner bias. In cases with non-determinism or noise, all-bias minimization, like all-variance minimization, becomes an approximation of the optimal approach.

The learning model discussed in this paper is a form of locally weighted regression (LWR) [Cleveland et al., 1988], which has been used in difficult machine learning tasks, notably the “robot juggler” of Schaal and Atkeson [1994]. Previous work [Cohn et al., 1995] discussed all-variance query selection for LWR; in the remainder of this paper, I describe a method for performing all-bias query selection. Section 2 describes the criterion that must be optimized for all-bias query selection. Section 3 describes the locally weighted regression learner used in this paper and describes how the all-bias criterion may be computed for it. Section 4 describes the results of experiments using this criterion on several simple domains. Directions for future work are discussed in Section 5.

2 All-bias query selection

Let us assume for the moment that we have a source of noise-free examples (x_i, y_i) and a deterministic learner which, given input x , outputs estimate $\hat{y}(x)$.² Let us also assume that we have an accurate estimate of the bias of \hat{y} which can be used to estimate the true function $y(x) = \hat{y}(x) - \text{bias}(x)$. We will break these rather strong assumptions of noise-free examples and accurate bias estimates in Section 4, but they are useful for deriving the theoretical approach described below.

²For clarity, I will drop the argument x except where required for disambiguation. I will also denote only the univariate case; the results apply in higher dimensions as well.

Given the accurate bias estimate, our task is then to force the biased estimator into the best approximation of $y(x)$ with the fewest number of examples. This, in effect, transforms the query selection problem into an example filter problem similar to that studied by Plutowski and White [1993] for neural networks. Below, I derive this criterion for estimating the change in error at x given a new queried example at \tilde{x} .

Since we have (temporarily) assumed a deterministic learner and noise-free data, the expected error in Equation 2 simplifies to:

$$E \left[(\hat{y}(x; \mathcal{D}) - y(x))^2 | x, \mathcal{D} \right] = (\hat{y}(x; \mathcal{D}) - y(x))^2 \quad (4)$$

We want to select a new \tilde{x} such that when we add (\tilde{x}, \tilde{y}) , the resulting squared bias is minimized:

$$(\hat{y}' - y)^2 \equiv (\hat{y}(x; \mathcal{D} \cup (\tilde{x}, \tilde{y})) - y(x))^2. \quad (5)$$

We will, for the remainder of the paper, use the “” to indicate estimates based on the initial training set plus the additional example (\tilde{x}, \tilde{y}) . To minimize Expression 5, we need to compute how a query at \tilde{x} will change the learner’s bias at x . If we assume that we know the input distribution,³ then we can integrate this change over the entire domain (using Monte Carlo procedures) to estimate the resulting average change, and select a \tilde{x} such that the expected squared bias is minimized. Defining $bias \equiv \hat{y} - y$ and $\Delta\hat{y} \equiv \hat{y}' - \hat{y}$, we can write the new squared bias as:

$$\begin{aligned} bias' &= (\hat{y}' - y)^2 \\ &= (\hat{y} + \Delta\hat{y} - y)^2 \\ &= \Delta\hat{y}^2 + 2\Delta\hat{y} \cdot bias + bias^2 \end{aligned} \quad (6)$$

Note that since $bias$ as defined here is independent of \tilde{x} , minimizing the bias is equivalent to minimizing $\Delta\hat{y}^2 + 2\Delta\hat{y} \cdot bias$.

The estimate of $bias'$ tells us how much our bias will change for a given \tilde{x} . We may optimize this value over \tilde{x} in one of a number of ways. In low dimensional spaces, it is often sufficient to consider a set of “candidate” \tilde{x} and select the one promising the smallest resulting error. In higher dimensional spaces, it is often more efficient to search for an optimal \tilde{x} with a response surface technique [Box and Draper, 1987], or hillclimb on $\partial bias' / \partial \tilde{x}$.

Estimates of $bias$ and $\Delta\hat{y}$ depend on the specific learning model being used. In Section 3, I describe a locally weighted regression model, and show how differentiable estimates of $bias$ and $\Delta\hat{y}$ may be computed for it.

2.1 An aside: why not just use $\hat{y} - \widehat{bias}$?

If we have an accurate bias estimate, it is reasonable to ask why we do not simply use the corrected $\hat{y} - \widehat{bias}$ as our predictor. Certainly, in the limit of a perfect bias estimate, the composite prediction would have zero bias,

³This assumption is contrary to the assumption normally made in some forms of learning, e.g. PAC-learning, but it is appropriate in many domains. If, for example, we are learning to control a robot arm, it is reasonable to assume that we know the distribution of positions over which we are interested in controlling it.

and we could concentrate solely on variance, as previous work has.

The answer to this question has several parts, the first of which is that for most learners, there are no perfect bias estimators. Bias estimators introduce their own bias and variance, which must be addressed in data selection.

We can define a composite learner which produces estimate $\hat{y}_c \equiv \hat{y} - \widehat{bias}$. Given a random training sample then, we would expect \hat{y}_c to outperform \hat{y} . However, there is no obvious way to select data for this composite learner other than selecting to maximize the performance of its two components. In our case, the second component (the bias estimate) is non-analytic, which leaves us selecting data so as to maximize the performance of the first component (the uncorrected estimator). We are now back to our original problem: we can select data so as to minimize either the bias or variance of the uncorrected LWR-based learner. Since the purpose of the correction is to give an unbiased estimator, intuition suggests that variance minimization would be the more sensible route in this case.

Regardless of how we select our data, we can use the composite estimator to make our predictions; depending on how noisy the bias estimate is, this may or may not improve the learner’s net performance. In the domains considered in this paper, I found that the performance of \hat{y}_c using random selection or variance minimization was not substantially different from that of the uncorrected \hat{y} (see Figure 7 in Section 4).

3 Locally weighted regression

The type of learner I consider here is a form of locally weighted regression (LWR) that is a slight variation on the LOESS model of Cleveland et al. [1988]. The LOESS model performs a linear regression on points in the data set, weighted by a kernel centered at x (see Figure 1). The kernel shape is a design parameter: the original LOESS model uses a “tricubic” kernel; in my experiments I use the more common Gaussian

$$h_i(x) \equiv h(x - x_i) = \exp(-k(x - x_i)^2),$$

where k is a smoothing parameter. For brevity, I will drop the argument x for $h_i(x)$, and define $n = \sum_i h_i$. We can then write the weighted means and covariances as:

$$\begin{aligned} \mu_x &= \sum_i h_i \frac{x_i}{n}, & \sigma_x^2 &= \sum_i h_i \frac{(x_i - x)^2}{n}, \\ \mu_y &= \sum_i h_i \frac{y_i}{n}, & \sigma_y^2 &= \sum_i h_i \frac{(y_i - \mu_y)^2}{n}, \\ \sigma_{y|x}^2 &= \sigma_y^2 - \frac{\sigma_{xy}^2}{\sigma_x^2}, & \sigma_{xy} &= \sum_i h_i \frac{(x_i - x)(y_i - \mu_y)}{n}. \end{aligned}$$

We use these means and covariances to produce an estimate \hat{y} at the x around which the kernel is centered, with a confidence term in the form of a variance estimate:

$$\begin{aligned} \hat{y} &= \mu_y + \frac{\sigma_{xy}}{\sigma_x^2}(x - \mu_x) \\ \sigma_{\hat{y}}^2 &= \sigma_{y|x}^2 \frac{\sum_i h_i^2}{n^2} \left[1 + \frac{(x - \mu_x)(x_i - \mu_x)}{\sigma_x^2} \right]^2. \end{aligned}$$

In all the experiments discussed in this paper, the smoothing parameter k was set so as to minimize $\sigma_{\hat{y}}^2$.

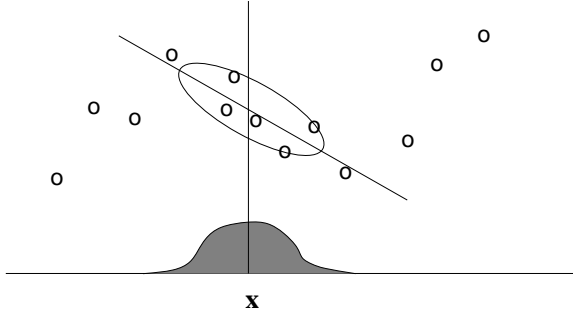


Figure 1: Locally weighted regression places a kernel around the point of interest x . The kernel is used to assign weightings to points in the training set, from which $\hat{y}(x)$ is computed via linear regression.

The low cost of incorporating new training examples makes this form of locally weighted regression appealing for learning systems which must operate in real time, or with time-varying target functions (e.g. [Schaal and Atkeson 1994]).

3.1 Computing $\Delta\hat{y}$ for LWR

If we know what new point (\tilde{x}, \tilde{y}) we're going to add, computing $\Delta\hat{y}$ for LWR is straightforward. Defining \tilde{h} as the weight given to \tilde{x} , we can write

$$\Delta\hat{y} = \hat{y}' - \hat{y} \quad (7)$$

$$\begin{aligned} &= \mu'_y + \frac{\sigma'_{xy}}{\sigma'^2_x}(x - \mu'_x) - \mu_y - \frac{\sigma_{xy}}{\sigma^2_x}(x - \mu_x) \\ &= h \frac{(\tilde{y} - \mu_y)}{n + h} - \frac{\sigma_{xy}}{\sigma^2_x}(x - \mu_x) \\ &\quad + \left(x - \frac{n\mu_x}{n + h} - \frac{h\tilde{x}}{n + h} \right) \\ &\quad \cdot \frac{(n + h)\sigma_{xy} + h \cdot (\tilde{x} - \mu_x)(\tilde{y} - \mu_y)}{(n + h)\sigma^2_x + h \cdot (\tilde{x} - \mu_x)^2} \end{aligned} \quad (8)$$

Note that computing $\Delta\hat{y}$ requires us to know both the \tilde{x} and \tilde{y} of the new point. In practice, we only know \tilde{x} . If we assume, however, that we can estimate the learner's bias at any x , then we can also estimate the unknown value $\tilde{y} \approx \hat{y}(\tilde{x}) - \text{bias}(\tilde{x})$. Below, I consider how to compute the bias estimate.

3.2 Estimating bias for LWR

The most common technique for estimating bias is cross-validation. Standard cross-validation however, only gives estimates of the bias at our specific training points, which are usually combined to form an average bias estimate. This is sufficient if one assumes that the training distribution is representative of the test distribution (which it isn't in query learning) and if one is content to just estimate the bias where one already has training data (which we can't be).

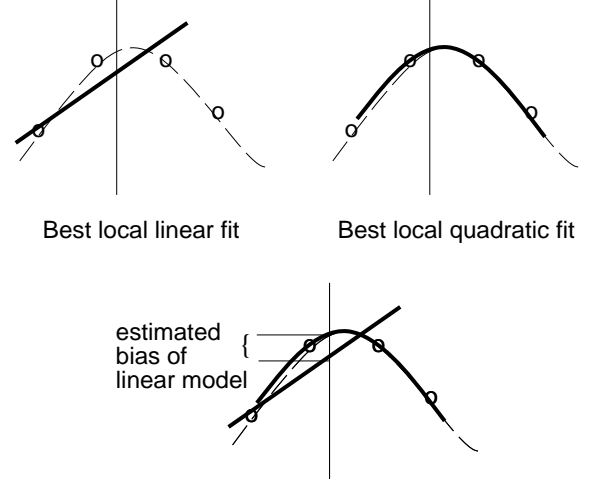


Figure 2: Box and Draper's method of estimating bias measures the difference between the estimator in question and a one-higher-order estimate.

In the query selection problem, we must be able to estimate the bias at all possible x . There are several ways we can get this estimate using LWR. Box and Draper [1987] suggest fitting a higher order model and measuring the difference. In the case of (linear) locally weighted regression, one would fit a locally quadratic regressor to the data and use the difference in estimates as the bias (see Figure 2). Under certain conditions on the higher-order bias terms, one can make some guarantees on the accuracy of this bias estimate.

The disadvantages of this method stem from the fact that it requires a higher order model. This requires additional computation to fit, and the fit is more prone to variance problems. For the experiments described in this paper, this method of bias estimation yielded poor results; two other bias-estimation techniques, however, performed very well.

3.2.1 Estimating bias by bootstrapping residuals

Another method of estimating bias is by bootstrapping the residuals of the training points. Based on the m available training points, and the predictor's fit to these points, a "bootstrap sample" is created by randomly drawing m values with replacement from the learner's residuals. These values are added to the original predictions to create a synthetic training set on which the learner is retrained.

By creating a number of bootstrapped predictions and comparing their average prediction with that of the original predictor, one arrives at a first-order bootstrap estimate of the predictor's bias [Connor 1993; Efron and Tibshirani, 1993]. It is known that this estimate is itself biased towards zero; a standard heuristic is to divide the estimate by 0.632 [Efron, 1983]. A disadvantage of the bootstrap method is that, because it requires repeated fitting, it is computationally expensive.

3.2.2 Estimating bias by fitting cross-validated estimates

One may also estimate the bias of a learner by fitting its own cross-validated residuals. We first compute the cross-validated residuals on the training examples. These produce estimates of the learner’s bias at each of the training points. We can then use these residuals as training examples for another learner (again LWR) to produce estimates of what the cross-validated error would be in places where we don’t have training data (see Figure 3).⁴

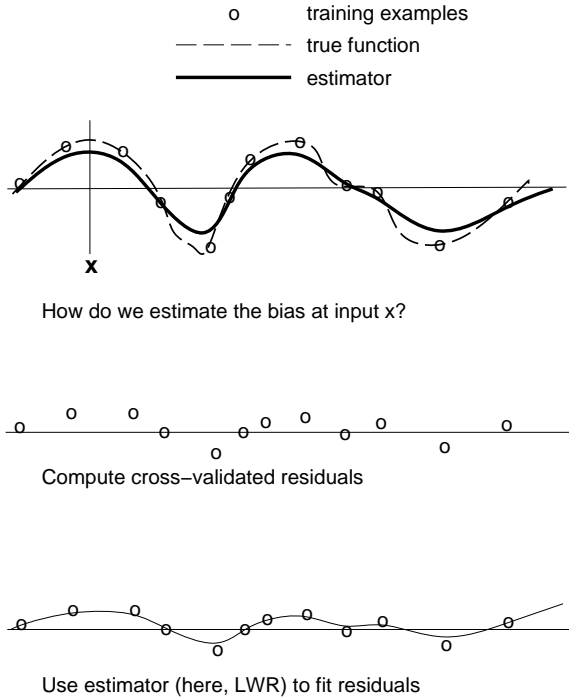


Figure 3: Applying the learner to its own cross-validated residuals produces an estimate of the bias that may be evaluated over the entire domain.

4 Empirical results

In the previous two sections, I have explained how having an estimate of $\Delta\hat{y}$ and *bias* for a learner allows one to compute the learner’s change in bias given a new query, and have shown how these estimates may be computed for a learner that uses locally weighted regression. Here, I apply these results to several simple problems using the “Arm2D” domain (Figure 4) and demonstrate that they may actually be used to select queries that minimize the statistical bias (and the error) of the learner.

⁴One subtlety that needs to be addressed is which residual is actually fit. Denote the cross-validated estimate as \hat{y}_{cv} . If we believe the data is noise-free, then the true value of the function at x is y , so the cross-validated bias is $\hat{y}_{cv} - y$. If, however, there is noise, we should assume that some of that misfit is due to noise. In this case, the proper bias estimate should be $\hat{y}_{cv} - \hat{y}$, with the remaining difference $\hat{y} - y$ being due to noise.

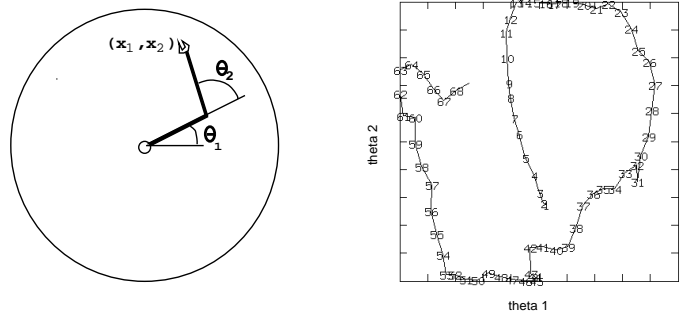


Figure 4: (left) Arm2D – The system learns arm kinematics by specifying joint angles (θ_1, θ_2) and observing tip coordinates (x_1, x_2) . The goal is to minimize the MSE of the learner’s model over the input distribution $(\Theta_1, \Theta_2) = (U[0, 2\pi], U[0, \pi])$. (right) A sample exploration trajectory in joint-space for the constrained arm problem, exploring according to the cross-validation-based bias minimizing criterion.

4.1 Bias estimates

I tested the accuracy of the three bias estimators by observing their correlations on 64 reference inputs, given 100 random training examples from the Arm2D domain. When corrected with the 632 heuristic described above, both the bootstrap and cross-validation methods produce fairly accurate, albeit noisy, bias estimates (Figure 5). The quadratic method produced poor correlation and was dropped from the study.

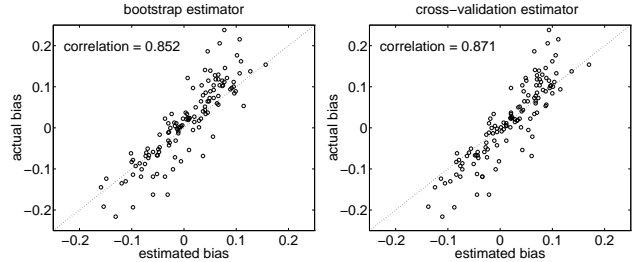


Figure 5: Correlations between estimated and actual biases for different estimators.

4.2 Bias minimization

I ran two series of experiments using the bias-minimizing criterion in conjunction with the bias estimation technique of the previous section on the “Arm2D” domain. The bias minimization criterion was used as follows: At each time step, the learner was given a set of 64 randomly chosen candidate queries and 64 uniformly chosen reference points. It evaluated $E'(x)$ for each reference point given each candidate point and selected for its next query the candidate point with the smallest average $E'(x)$ over the reference points. I compared the bias-minimizing strategy (using the cross-validation and bootstrap estimation techniques) against random sampling and the variance-minimizing strategy discussed in Cohn et al. [1995]. On a Sparc 10, with m training examples, the average evaluation times per candidate per

reference point were $58 + 0.16m$ μ seconds for the variance criterion, $65 + 0.53m$ μ seconds for the cross-validation-based bias criterion, and $83 + 3.7m$ μ seconds for the bootstrap-based bias criterion (with 20x resampling).

To test whether the bias-only assumption was robust against the presence of noise, 1% Gaussian noise was added to the input values of the training data in all experiments. This simulates noisy position effectors on the arm, and results in non-Gaussian noise in the output coordinate system.

In the first series of experiments, the candidate points were drawn uniformly over $(U[0, 2\pi], U[0, \pi])$. In unconstrained domains like this, random sampling is a fairly good default strategy. The bias minimization strategies still significantly outperform both random sampling and the variance minimizing strategy in these experiments (see Figure 6).

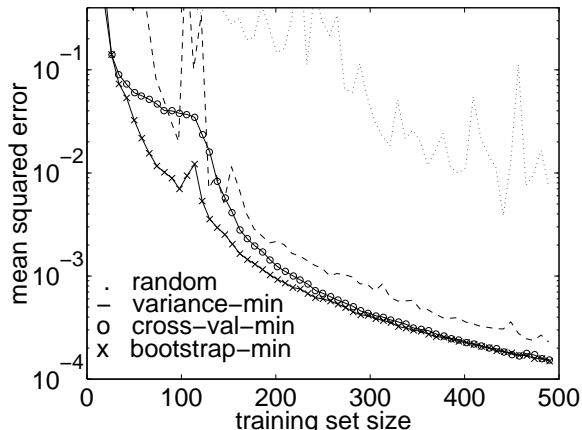


Figure 6: MSE as a function of number of noisy training examples for the unconstrained arm problem. The cross-validation and bootstrap bias-minimization strategies give a factor of 3 improvement over random selection, and a slight improvement over variance-only minimization. Errors are averaged over 10 runs for the bootstrap method and 15 runs for all others. One run with the cross-validation-based method was excluded when k failed to converge to a reasonable value.

In the second series of experiments, candidates were drawn uniformly from a region local to the previously selected query: $(\hat{\theta}_1 \pm 0.2\pi, \hat{\theta}_2 \pm 0.1\pi)$. This corresponds to restricting the arm to local motions. In a constrained problem such as this, random sampling is a poor strategy; both the bias and variance-reducing strategies outperform it at least an order of magnitude. Further, the bias-minimization strategy outperforms variance minimization by a large margin (Figure 7). Figure 4 shows an exploration trajectory produced by pursuing the bias-minimizing criterion. It is noteworthy that, although the implementation in this case was a greedy (one-step) minimization, the trajectory results in globally good exploration.

5 Discussion

I have argued in this paper that, in many situations, selecting queries to minimize learner bias is an appropriate

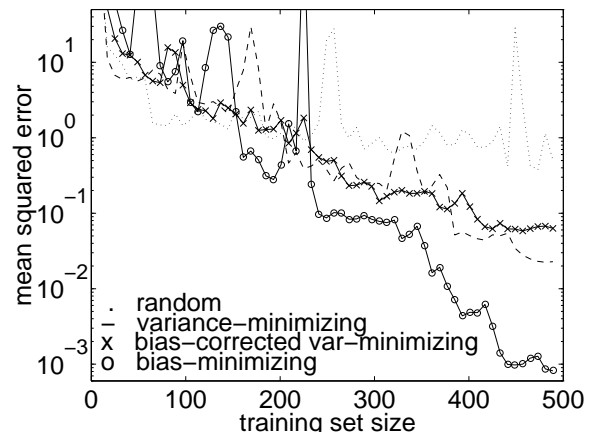


Figure 7: MSE as a function of number of noisy training examples for the constrained arm problem. Bias-minimization significantly outperforms the variance-minimizing algorithm and random exploration.

and effective strategy for active learning. I have given empirical evidence that, with a LWR-based learner and the examples considered here, the strategy is effective even in the presence of noise.

Beyond minimizing either bias or variance, an important next step is to explicitly minimize them together. The bootstrap-based estimate should facilitate this, as it produces a complementary variance estimate with little additional computation.⁵ By optimizing over both criteria simultaneously, we expect to derive a criterion that that, in terms of statistics, is truly optimal for selecting queries.

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⁵Because it assumes that all misfit is due to variance, the variance estimate in Cohn et al. [1995] may be inaccurate if significant bias is present.

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